

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: ssspta1653adk

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS 1	Web Page URLs for STN Seminar Schedule - N. America
NEWS 2	"Ask CAS" for self-help around the clock
NEWS 3 SEP 01	New pricing for the Save Answers for SciFinder
Wizard within	
NEWS 4 OCT 28	STN Express with Discover!
NEWS 5 NOV 30	KOREAPAT now available on STN
NEWS 6 DEC 01	PHAR reloaded with additional data
NEWS 7 DEC 09	LISA now available on STN
2004	12 databases to be removed from STN on December 31, 2004
NEWS 8 DEC 15	MEDLINE update schedule for December 2004
NEWS 9 DEC 17	ELCOM reloaded; updating to resume; current-awareness
NEWS 10 DEC 17	alerts (SDIs) affected
awareness	COMPUAB reloaded; updating to resume; current-
NEWS 11 DEC 17	alerts (SDIs) affected
awareness	SOLIDSTATE reloaded; updating to resume; current-
NEWS 12 DEC 17	alerts (SDIs) affected
awareness	CERAB reloaded; updating to resume; current-
NEWS 13 DEC 17	alerts (SDIs) affected
	THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB
NEWS EXPRESS	OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer

agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 10:14:29 ON 28 DEC 2004

=> fil reg
COST IN U.S. DOLLARS SINCE FILE
TOTAL ENTRY
SESSION
FULL ESTIMATED COST 0.21
0.21

FILE 'REGISTRY' ENTERED AT 10:14:35 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9
DICTIONARY FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9

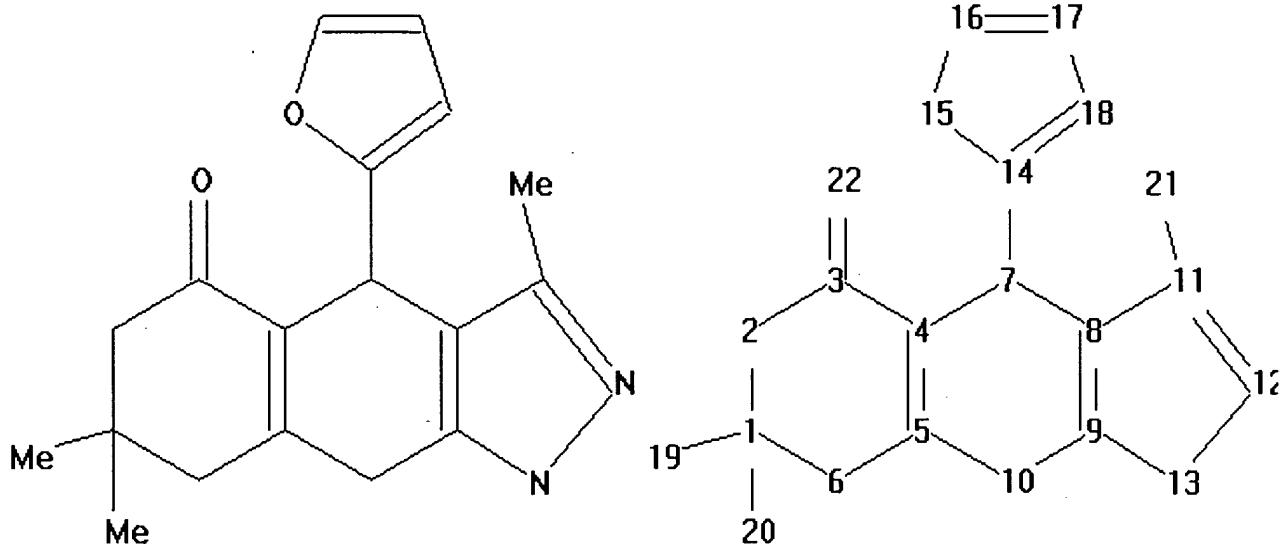
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> Uploading H:\STN queries\10612885.str



chain nodes :

19 20 21 22

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18

chain bonds :

1-19 1-20 3-22 7-14 11-21

ring bonds :

1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18

exact/norm bonds :

1-2 1-6 2-3 3-4 3-22 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10
9-13 11-12 12-13 14-15 14-18 15-16 16-17 17-18

exact bonds :

1-19 1-20 7-14 11-21

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS

L1 STRUCTURE UPLOADED

=> S 11

SAMPLE SEARCH INITIATED 10:14:50 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 8 TO ITERATE

100.0% PROCESSED 8 ITERATIONS 0

ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 8 TO 329

PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 ful
FULL SEARCH INITIATED 10:14:53 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 166 TO ITERATE

100.0% PROCESSED 166 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

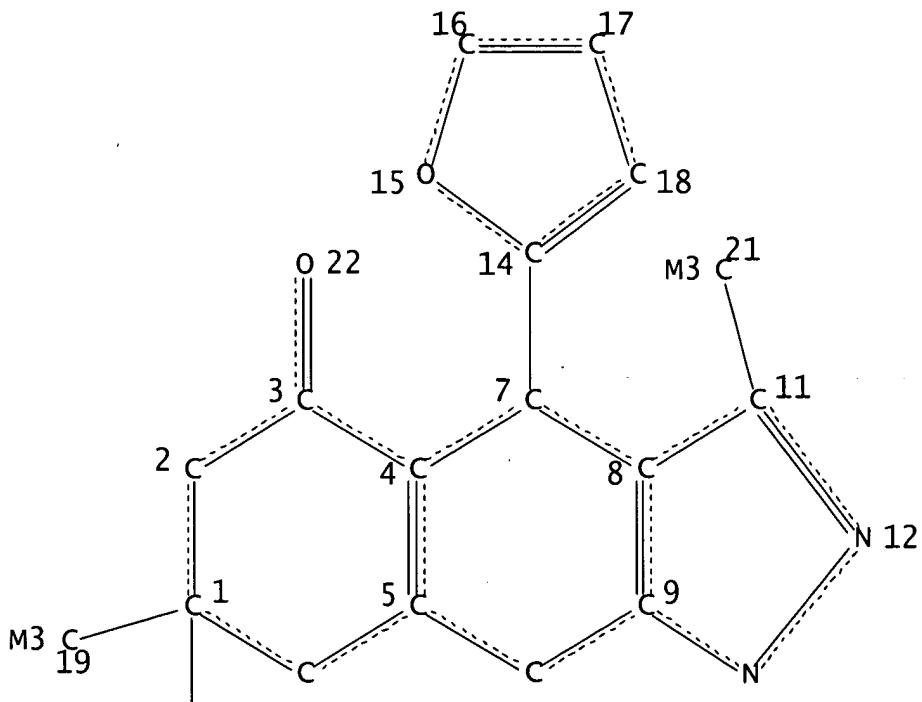
L3 0 SEA SSS FUL L1

=> s l1 ful fam
FULL SEARCH INITIATED 10:14:58 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

L4 0 SEA FAM FUL L1

=> d l1
L1 HAS NO ANSWERS
L1 STR



Page 1-A

6 10 13
C M3
20

Page 2-A

NODE ATTRIBUTES:

HCOUNT IS M3 AT 19
HCOUNT IS M3 AT 20
HCOUNT IS M3 AT 21
NSPEC IS R AT 1
NSPEC IS R AT 2
NSPEC IS R AT 3
NSPEC IS R AT 4
NSPEC IS R AT 5
NSPEC IS R AT 6
NSPEC IS R AT 7
NSPEC IS R AT 8
NSPEC IS R AT 9
NSPEC IS R AT 10
NSPEC IS R AT 11
NSPEC IS R AT 12
NSPEC IS R AT 13
NSPEC IS R AT 14
NSPEC IS R AT 15
NSPEC IS R AT 16
NSPEC IS R AT 17
NSPEC IS R AT 18
NSPEC IS C AT 19
NSPEC IS C AT 20
NSPEC IS C AT 21
NSPEC IS C AT 22

DEFAULT MLEVEL IS ATOM

MLEVEL IS CLASS AT 19 20 21 22

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

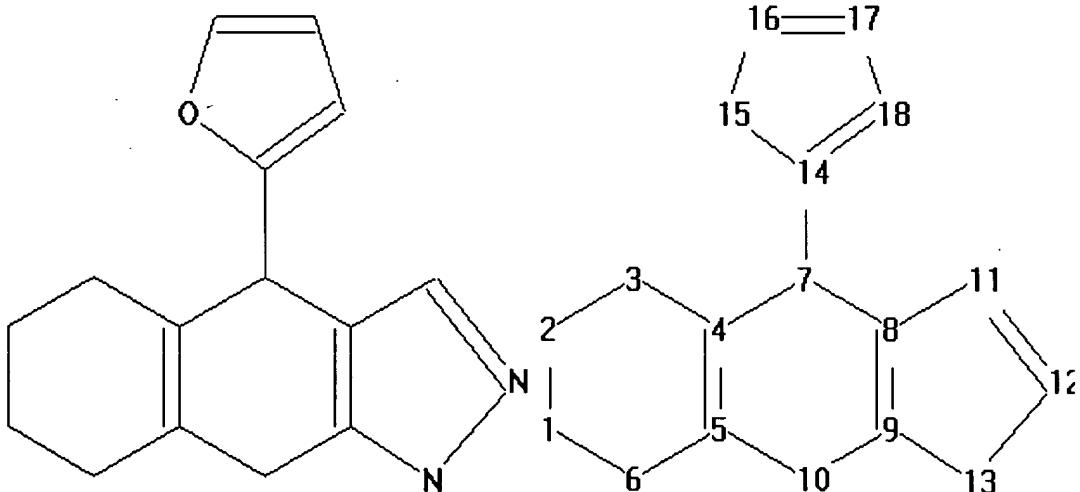
RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

=>

Uploading H:\STN queries\10612885a.str



ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
7-14
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
7-14

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom

L5 STRUCTURE UPLOADED

=> s 15
SAMPLE SEARCH INITIATED 10:18:34 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 24 TO ITERATE

100.0% PROCESSED 24 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
PROJECTED ITERATIONS: 187 TO 773
BATCH **COMPLETE**
PROJECTED ANSWERS: 0 TO 0

L6 0 SEA SSS SAM L5

=> s 15 ful
FULL SEARCH INITIATED 10:18:38 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 607 TO ITERATE

100.0% PROCESSED 607 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

L7 0 SEA SSS FUL L5

=> s 15 fam
SAMPLE SEARCH INITIATED 10:18:42 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L8 0 SEA FAM SAM L5

=> S 15 fam ful
FULL SEARCH INITIATED 10:18:48 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

L9 0 SEA FAM FUL L5

=> DIS HIST

(FILE 'HOME' ENTERED AT 10:14:29 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:14:35 ON 28 DEC 2004

L1 STRUCTURE uploaded
L2 0 S L1
L3 0 S L1 FUL
L4 0 S L1 FUL FAM
L5 STRUCTURE uploaded
L6 0 S L5
L7 0 S L5 FUL
L8 0 S L5 FAM
L9 0 S L5 FAM FUL

=>

---Logging off of STN---

=>
Executing the logoff script...

=> LOG Y

COST IN U.S. DOLLARS	SINCE FILE
TOTAL	ENTRY
SESSION	
FULL ESTIMATED COST	434.86
435.07	

STN INTERNATIONAL LOGOFF AT 10:19:23 ON 28 DEC 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1653adk

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

***** Welcome to STN International *****

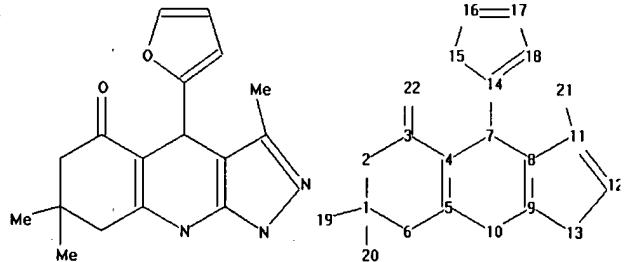
NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 SEP 01 New pricing for the Save Answers for SciFinder Wizard within
Wizard within STN Express with Discover!
NEWS 4 OCT 28 KOREAPAT now available on STN
NEWS 5 NOV 30 PHAR reloaded with additional data
NEWS 6 DEC 01 LISA now available on STN
NEWS 7 DEC 09 12 databases to be removed from STN on December 31, 2004
NEWS 8 DEC 15 MEDLINE update schedule for December 2004
NEWS 9 DEC 17 ELCOM reloaded; updating to resume; current-awareness
NEWS 10 DEC 17 alerts (SDIs) affected
COMPUAB reloaded; updating to resume; current-awareness
NEWS 11 DEC 17 alerts (SDIs) affected
SOLIDSTATE reloaded; updating to resume; current-awareness
NEWS 12 DEC 17 alerts (SDIs) affected
CERAB reloaded; updating to resume; current-awareness
NEWS 13 DEC 17 alerts (SDIs) affected
THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB

NEWS EXPRESS OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT MACINTOSH VERSION IS V6.0C(ENG) AND V6.0C(JP)
AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to
STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer



```

chain nodes :
19 20 21 22
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
1-19 1-20 3-22 7-14 11-21
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 3-22 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10
9-13 11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
1-19 1-20 7-14 11-21

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS

L1 STRUCTURE UPLOADED

=> d 11
L1 HAS NO ANSWERS
L1 STR

agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

ETLE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004

=> FIL REGISTRY
COST IN U.S. DOLLARS
TOTAL
SESSION
FULL ESTIMATED COST
0.21
SINCE FILE
ENTRY
0.21

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9
DICTIONARY FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004
Please note that search-term pricing does apply when

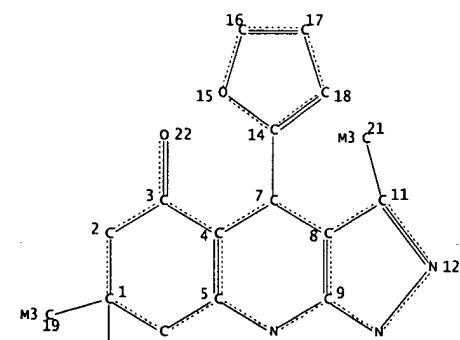
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. P

information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> Uploading H:\STN queries\10612885b.str



Page 1-A

20 C M3 6 10 13

Page 2-A		NODE ATTRIBUTES:	
HCOUNT	IS M3	AT	19
HCOUNT	IS M3	AT	20
HCOUNT	IS M3	AT	21
NSPEC	IS R	AT	1
NSPEC	IS R	AT	2
NSPEC	IS R	AT	3
NSPEC	IS R	AT	4
NSPEC	IS R	AT	5
NSPEC	IS R	AT	6
NSPEC	IS R	AT	7
NSPEC	IS R	AT	8
NSPEC	IS R	AT	9
NSPEC	IS R	AT	10
NSPEC	IS R	AT	11
NSPEC	IS R	AT	12
NSPEC	IS R	AT	13
NSPEC	IS R	AT	14
NSPEC	IS R	AT	15
NSPEC	IS R	AT	16
NSPEC	IS R	AT	17
NSPEC	IS R	AT	18
NSPEC	IS C	AT	19
NSPEC	IS C	AT	20

NSPEC IS C AT 21
 NSPEC IS C AT 22
 DEFAULT MLEVEL IS ATOM
 MLEVEL IS CLASS AT 19 20 21 22
 DEFAULT ELEVEL IS LIMITED

GRAPH ATTRIBUTES:
 RING(S) ARE ISOLATED OR EMBEDDED
 NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

=> s 11 sam
 SAMPLE SEARCH INITIATED 10:25:56 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS
 ANSWERS
 SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
 PROJECTED ITERATIONS: 0 TO 0
 PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s 11 fam
 SAMPLE SEARCH INITIATED 10:26:03 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS
 ANSWERS
 SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
 PROJECTED ITERATIONS: 0 TO 0
 PROJECTED ANSWERS: 0 TO 0

L3 0 SEA FAM SAM L1

=> s 11 ful
 FULL SEARCH INITIATED 10:26:08 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS
 ANSWERS
 SEARCH TIME: 00.00.01

L4 8 SEA SSS FUL L1

=> s 11 ful fam
 FULL SEARCH INITIATED 10:26:12 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 5 TO ITERATE

100.0% PROCESSED 5 ITERATIONS

ANSWERS
 SEARCH TIME: 00.00.01

L5 1 SEA FAM FUL L1

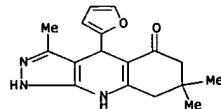
=> 15 not 14
 L6 0 LS NOT L4

=> d 15 1 all

LS ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 645337-25-3 REGISTRY
 ED Entered STN: 03 Feb 2004
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C17 H19 N3 O2
 SR CA
 LC STN Files: CA, CAPLUS, USPATFULL
 DT CA Caplus document type: Patent
 RL.P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Count
C40	O4	5	C40	16.138.5	1
C3N2-CSN-C6 N2C3-NCS-C6	5-6-6		C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1	pH 1	(1) ACD
Bioconc. Factor (BCF)	8.39	pH 4	(1) ACD
Bioconc. Factor (BCF)	197.8	pH 7	(1) ACD

Bioconc. Factor (BCF) 98.8 | pH 8 | (1) ACD
 Bioconc. Factor (BCF) 98.8 | pH 10 | (1) ACD
 Boiling Point (BP) 478.2+-45.0 deg C | 760 Torr | (1) ACD
 Enthalpy of Vap. (HVAP) 74.23+-3.0 kJ/mol | (1) ACD
 Flash Point (FP) 243.0+-51.7 deg C | (1) ACD
 Freely Rotatable Bonds (FRB) 1 | (1) ACD
 H acceptors (HAC) 5 | (1) ACD
 H donors (HD) 2 | (1) ACD
 Koc (KOC) 1 | pH 1 | (1) ACD
 Koc (KOC) 79.0 | pH 4 | (1) ACD
 Koc (KOC) 922 | pH 7 | (1) ACD
 Koc (KOC) 931 | pH 8 | (1) ACD
 Koc (KOC) 932 | pH 10 | (1) ACD
 LogD (LOGD) -0.49 | pH 1 | (1) ACD
 LogD (LOGD) 1.86 | pH 4 | (1) ACD
 LogD (LOGD) 2.92 | pH 7 | (1) ACD
 LogD (LOGD) 2.93 | pH 8 | (1) ACD
 LogD (LOGD) 2.93 | pH 10 | (1) ACD
 LogP (LOGP) 2.928+-0.412 | (1) ACD
 Molar Solubility (SLB.MOL) >0.1 - <1 mol/L | pH 1 | (1) ACD
 Molar Solubility (SLB.MOL) <0.01 mol/L | pH 4 | (1) ACD
 Molar Solubility (SLB.MOL) <0.01 mol/L | pH 7 | (1) ACD
 Molar Solubility (SLB.MOL) <0.01 mol/L | pH 8 | (1) ACD
 Molar Solubility (SLB.MOL) <0.01 mol/L | pH 10 | (1) ACD
 Molecular Weight (MW) 297.35 | (1) ACD
 pKa (PKA) 13.94+-0.60 | Most Acidic | (1) ACD
 pKa (PKA) 4.82+-0.60 | Most Basic | (1) ACD
 Vapor Pressure (VP) 2.62E-09 Torr | 25 deg C | (1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs)
 Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, US 2004171541 A1 20040902 US 2003-613754 20030702 US 2004116346 A1 20040617 US 2003-612885 20030703 PRAI US 2002-393360P 20020703 US 2002-393361P 20020703 US 2002-394110P 20020703 AB Compds. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazolopyrimidine, pyridazinone, pyridine and piperidine.

ST EPO receptor modulator small mol
 IT Proteins
 RL: BSU (Biological study, unclassified); BIOL (Biological study)
 (Bcl-xL, expression; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
 IT Peptides, biological studies
 RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (biological study); USES (Uses)
 (EPO receptor modulating sequence; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
 IT Cell membrane
 (EPO receptors of; affinity small mols. for erythropoietin

See HELP PROPERTIES for information about property data sources in REGISTRY.
 1 REFERENCES IN FILE CA (1907 TO DATE)
 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 140:87744 CA Full-text
 TI Affinity small molecules for the EPO receptor
 IN Olsson, Lennart; Naranda, Tatjana
 PA Receptor, Inc., USA
 SO PCT Int. Appl., 85 pp.
 CODEN: PIXXD2

DT Patent

LA English

IC ICM C07K

CC 1-12 (Pharmacology)

Section cross-reference(s): 2

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO. DATE
 PI WO 2004005323 A2 20040115 WO 2003-US21394 20030703
 WO 2004005323 A3 20040701

(EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Anemia (disease)
Cell proliferation
Combinatorial library
Drug delivery systems
Drug screening
Erythrocyte
Erythropoiesis
Hematocrit
Hematopoietic precursor cell
Human
Reticulocyte
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Nerve
(neuron; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Cytoprotective agents
(neuroprotective; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 2503-56-2 40775-78-8 S1646-16-3 S1646-17-4 S1646-19-6
S1646-43-6 56347-20-7 63901-48-4 90559-98-1 90815-61-5 113967-71-8
113967-74-1 194342-06-8 212074-47-0 244167-89-3 245082-87-5
245413-82-5 259683-29-9 261704-08-9 261704-09-0 262291-81-6

16-1 263267-38-5 287728-46-5 303145-64-4 303145-73-5 338793-645337-19-5 645337-20-8 645337-21-9 645337-22-0 645337-23-1 645337-24-2 645337-25-3
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
EPO (affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 11096-26-7, Erythropoietin
RL: BSU (Biological study, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 239133-03-0 645415-22-1
RL: PRP (Properties)
(unclaimed sequence; affinity small mols. for the EPO receptor)

=> d 14 1-8 ful
'FUL' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN
CALC - Table of calculated properties
EPROP - Table of experimental properties
PROP - EPROP and CALC

Any CA File format may be combined with any substance format to

obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL
IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OBIBB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.

ENTER DISPLAY FORMAT (IDE):all

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

RN 748146-78-3 REGISTRY

ED Entered STN: 20 Sep 2004

CH INDEX NAME NOT YET ASSIGNED

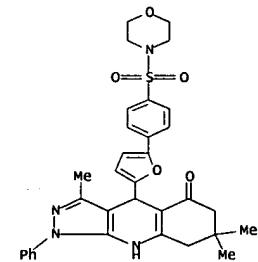
FS 3D CONCORD

MF C33 H34 N4 O5 S

SR Chemical Library

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	2
C4NO	NC20C2	6	C4NO	46.402.1	1
C3N2-CSN-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	2.13	pH 1	(1) ACD
Bioconc. Factor (BCF)	649	pH 4	(1) ACD
Bioconc. Factor (BCF)	1353	pH 7	(1) ACD
Bioconc. Factor (BCF)	1354	pH 8	(1) ACD
Bioconc. Factor (BCF)	1355	pH 10	(1) ACD
Boiling Point (BP)	749.7+-70.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	109.28+-3.0 kJ/mol		(1) ACD
Flash Point (FP)	407.2+-64.2 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	4		(1) ACD
H acceptors (HAC)	9		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	9.55	pH 1	(1) ACD
Koc (KOC)	2910	pH 4	(1) ACD
Koc (KOC)	6062	pH 7	(1) ACD
Koc (KOC)	6067	pH 8	(1) ACD
Koc (KOC)	6068	pH 10	(1) ACD
LogD (LOGD)	1.62	pH 1	(1) ACD
LogD (LOGD)	4.10	pH 4	(1) ACD
LogD (LOGD)	4.42	pH 7	(1) ACD
LogD (LOGD)	4.42	pH 8	(1) ACD
LogD (LOGD)	4.42	pH 10	(1) ACD
LogP (LOGP)	4.424+-0.655		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD

Molar Solubility (SLB.MOL) <0.01 mol/L | pH 10 | (1) ACD
 Molecular Weight (MW) 598.71 | (1) ACD
 pKa (PKA) 3.95+/-0.60 | Most Basic | (1) ACD
 Vapor Pressure (VP) 2.33E-22 Torr | 25 deg C | (1) ACD

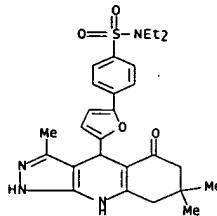
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748146-41-0 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C27 H32 N4 O4 S
 SR Chemical Library

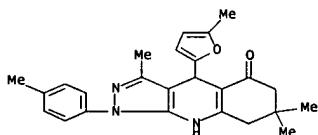
Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System RF	Ring Identifier RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	1
C3N2-C5N-C6 N2C3-NCS-C6 5-6-6			C10N3	1894.85.24	1



Calculated Properties (CALC)

C40 | OC4 | 5 | C40 | 16.138.5 | 1
 C6 | C6 | 6 | C6 | 46.150.18 | 1
 C3N2-C5N-C6|N2C3-NCS-C6|5-6-6 | C10N3 | 1894.85.24 | 1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	2.22	pH 1	(1) ACD
Bioconc. Factor (BCF)	847	pH 4	(1) ACD
Bioconc. Factor (BCF)	2188	pH 7	(1) ACD
Bioconc. Factor (BCF)	2191	pH 8	(1) ACD
Bioconc. Factor (BCF)	2192	pH 10	(1) ACD
Boiling Point (BP)	536.3+/-50.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	81.29+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	278.1+/-54.2 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	2		(1) ACD
H acceptors (HAC)	5		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	8.69	pH 1	(1) ACD
Koc (KOC)	3308	pH 4	(1) ACD
Koc (KOC)	8549	pH 7	(1) ACD
Koc (KOC)	8562	pH 8	(1) ACD
Koc (KOC)	8563	pH 10	(1) ACD
logD (LOGD)	1.70	pH 1	(1) ACD
logD (LOGD)	4.29	pH 4	(1) ACD
logD (LOGD)	4.70	pH 7	(1) ACD
logD (LOGD)	4.70	pH 8	(1) ACD
logD (LOGD)	4.70	pH 10	(1) ACD
logP (LOGP)	4.699+/-0.511		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	401.50		(1) ACD
pKa (PKA)	4.16+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	1.42E-11 Torr	25 deg C	(1) ACD

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1.22	pH 1	(1) ACD
Bioconc. Factor (BCF)	274	pH 4	(1) ACD
Bioconc. Factor (BCF)	3190	pH 7	(1) ACD
Bioconc. Factor (BCF)	3221	pH 8	(1) ACD
Bioconc. Factor (BCF)	3225	pH 10	(1) ACD
Boiling Point (BP)	680.7+/-65.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	99.88+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	365.5+/-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	5		(1) ACD
H acceptors (HAC)	8		(1) ACD
H donors (HD)	2		(1) ACD
Koc (KOC)	4.28	pH 1	(1) ACD
Koc (KOC)	958	pH 4	(1) ACD
Koc (KOC)	11170	pH 7	(1) ACD
Koc (KOC)	11278	pH 8	(1) ACD
Koc (KOC)	11289	pH 10	(1) ACD
logD (LOGD)	1.50	pH 1	(1) ACD
logD (LOGD)	3.85	pH 4	(1) ACD
logD (LOGD)	4.91	pH 7	(1) ACD
logD (LOGD)	4.92	pH 8	(1) ACD
logD (LOGD)	4.92	pH 10	(1) ACD
logP (LOGP)	4.920+/-0.532		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	508.63		(1) ACD
pKa (PKA)	13.94+/-0.60	Most Acidic	(1) ACD
pKa (PKA)	4.82+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	2.19E-18 Torr	25 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748145-15-5 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C25 H27 N3 O2
 SR Chemical Library

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System RF	Ring Identifier RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	1
C3N2-C5N-C6 N2C3-NCS-C6 5-6-6			C10N3	1894.85.24	1

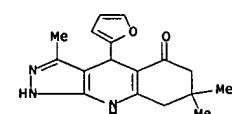
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 645337-25-3 REGISTRY
 ED Entered STN: 03 Feb 2004
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C17 H19 N3 O2
 SR CA
 LC STN Files: CA, CAPLUS, USPATFULL
 DT.CA Cplus document type: Patent
 RL.P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System RF	Ring Identifier RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C3N2-C5N-C6 N2C3-NCS-C6 5-6-6			C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1	pH 1	(1) ACD
Bioconc. Factor (BCF)	8.39	pH 4	(1) ACD
Bioconc. Factor (BCF)	97.8	pH 7	(1) ACD
Bioconc. Factor (BCF)	98.8	pH 8	(1) ACD
Bioconc. Factor (BCF)	98.8	pH 10	(1) ACD
Boiling Point (BP)	478.2+/-45.0 deg C	760 Torr	(1) ACD

Enthalpy of Vap. (HVAP)	74.23+/-3.0 kJ/mol	(1) ACD
Flash Point (FP)	243.0+/-51.7 deg C	(1) ACD
Freely Rotatable Bonds (FRB)	1	(1) ACD
H acceptors (HAC)	5	(1) ACD
H donors (HD)	2	(1) ACD
Koc (KOC)	1	pH 1
Koc (KOC)	79.0	pH 4
Koc (KOC)	922	pH 7
Koc (KOC)	931	pH 8
Koc (KOC)	932	pH 10
logP (LOGD)	-0.49	pH 1
logP (LOGD)	1.86	pH 4
logP (LOGD)	2.92	pH 7
logP (LOGD)	2.93	pH 8
logP (LOGD)	2.93	pH 10
logP (LOGP)	2.928+/-0.412	(1) ACD
Molar Solubility (SLB.MOL)	>0.1 - <1 mol/L	pH 1
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10
Molecular Weight (MW)	297.35	(1) ACD
pKa (PKA)	13.94+/-0.60	Most Acidic
pKa (PKA)	4.82+/-0.60	Most Basic
Vapor Pressure (VP)	2.62E-09 Torr	25 deg C
		(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN	140:87744	CA Full-text
TI	Affinity small molecules for the EPO receptor	
IN	Olsson, Lennart; Naranda, Tatjana	
PA	Reception, Inc., USA	
SO	PCT Int. Appl., 85 pp.	
	CODEN: PIXXD2	
DT	Patent	
LA	English	
IC	ICM C07K	
CC	1-12 (Pharmacology)	
	Section cross-reference(s): 2	
FAN.CNT 1		
	PATENT NO.	KIND
PI	WO 2004005323	A2
	WO 2004005323	A3
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
		CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,
		IT
		APPLICATION NO. DATE
WO 2004005323	20040115	WO 2003-US21394 20030703
WO 2004005323	20040701	

the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Anemia (disease)
Cell proliferation
Combinatorial library
Drug delivery systems
Drug screening
Erythrocyte
Erythropoiesis
Hematocrit
Hematopoietic precursor cell
Human
Reticulocyte
(affinity small mols. for erythropoietin (EPO) receptor and receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Nerve
(neuron; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Cytoprotective agents
(neuroprotective; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 2503-56-2 40775-78-8 51646-16-3 51646-17-4 51646-19-6
51646-43-6
56347-20-7 63901-48-4 90559-98-1 90815-61-5 113967-71-8
113967-74-1 194342-06-8 212074-47-0 244167-89-3 245082-87-5
245413-82-5 259683-29-9 261704-08-9 261704-09-0 262291-81-6
263267-38-5 287728-46-5 303145-64-4 303145-73-5 338793-16-1
645337-19-5 645337-20-8 645337-21-9 645337-22-0 645337-

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW, RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, US 2004171541 A1 20040902 US 2003-613754 20030702 US 2004116346 A1 20040617 US 2003-612885 20030703

PRAI US 2002-393360P 20020703
US 2002-393361P 20020703
US 2002-394110P 20020703

AB Compds. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazolopyrimidine, pyridazinone, pyridine and piperidine.

ST EPO receptor modulator small mol

IT Proteins
RL: BSU (Biological study, unclassified); BIOL (Biological study)
(Bcl-xL, expression; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Peptides, biological studies
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(EPO receptor modulating sequence; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Cell membrane
(EPO receptors of; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

23-1 645337-24-2 645337-25-3
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 11096-26-7 Erythropoietin
RL: BSU (Biological study, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 239133-03-0 645415-22-1
RL: PRP (Properties)
(unclaimed sequence; affinity small mols. for the EPO receptor)

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 521318-71-8 REGISTRY
ED Entered STN: 28 May 2003
CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-4-(5-methyl-2-furanyl)-(9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C24 H27 N5 O2
SR Chemical Library
LC STN Files: CHEMCATS

Ring System Data

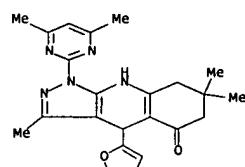
Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System Formula	Ring Identifier	Ring Occurrence Count
C40	O[OC4]	5	C40	16.138.5	1
C4N2	[NCN3]	6	C4N2	46.195.39	1
C3N2-C5N-C6[NC2-NC5-C6]5-6-6			C10N3	1894.85.241	1

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 521284-01-5 REGISTRY
 ED Entered STN: 28 May 2003
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyridinyl)-4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C23 H25 N5 O2
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System Formula	Ring Identifier	RID
EA	ES	SZ	RF	RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C4N2	[N]CNC3	6	C4N2	46.195.39	1
C3N2-C5N-C6	[N]2C3-NCS-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	242	pH 1	(1) ACD
Bioconc. Factor (BCF)	633	pH 4	(1) ACD
Bioconc. Factor (BCF)	634	pH 7	(1) ACD
Bioconc. Factor (BCF)	634	pH 8	(1) ACD
Bioconc. Factor (BCF)	634	pH 10	(1) ACD
Boiling Point (BP)	607.6+-65.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	90.29+-3.0 kJ/mol		(1) ACD
Flash Point (FP)	321.2+-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	2		(1) ACD
H acceptors (HAC)	7		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	1346	pH 1	(1) ACD
Koc (KOC)	3518	pH 4	(1) ACD
Koc (KOC)	3524	pH 7	(1) ACD
Koc (KOC)	3524	pH 8	(1) ACD
Koc (KOC)	3524	pH 10	(1) ACD
logD (LOGD)	3.57	pH 1	(1) ACD
logD (LOGD)	3.99	pH 4	(1) ACD
logD (LOGD)	3.99	pH 7	(1) ACD
logD (LOGD)	3.99	pH 8	(1) ACD
logD (LOGD)	3.99	pH 10	(1) ACD
logP (LOGP)	3.990+-0.891		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	417.50		(1) ACD
pKa (PKA)	1.15+-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	1.04E-14 Torr	25.0 deg C	(1) ACD

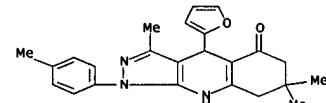
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software

Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	108	pH 1	(1) ACD
Bioconc. Factor (BCF)	283	pH 4	(1) ACD
Bioconc. Factor (BCF)	283	pH 7	(1) ACD

Bioconc. Factor (BCF)	283	pH 8	(1) ACD
Bioconc. Factor (BCF)	283	pH 10	(1) ACD
Boiling Point (BP)	601.5+-65.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	89.51+-3.0 kJ/mol		(1) ACD
Flash Point (FP)	317.6+-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	2		(1) ACD
H acceptors (HAC)	7		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	756	pH 1	(1) ACD
Koc (KOC)	1977	pH 4	(1) ACD
Koc (KOC)	1980	pH 7	(1) ACD
Koc (KOC)	1980	pH 8	(1) ACD
Koc (KOC)	1980	pH 10	(1) ACD
logD (LOGD)	3.11	pH 1	(1) ACD
logD (LOGD)	3.53	pH 4	(1) ACD
logD (LOGD)	3.53	pH 7	(1) ACD
logD (LOGD)	3.53	pH 8	(1) ACD
logD (LOGD)	3.53	pH 10	(1) ACD
logP (LOGP)	3.530+-0.890		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	403.48		(1) ACD
pKa (PKA)	1.15+-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	2.00E-14 Torr	25.0 deg C	(1) ACD

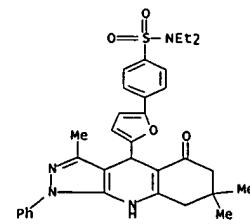
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software



L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 378189-53-8 REGISTRY
 ED Entered STN: 26 Dec 2001
 CN Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-1H-pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C33 H36 N4 O4 S
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System Formula	Ring Identifier	RID
EA	ES	SZ	RF	RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	2
C3N2-C5N-C6	[N]2C3-NCS-C6	5-6-6	C10N3	1894.85.24	1



Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System	Ring Identifier	RID
EA	ES	SZ	RF	RID	Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	2
C3N2-C5N-C6	[N]2C3-NCS-C6	5-6-6	C10N3	1894.85.24	1

Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	22.5	pH 1	(1) ACD
Bioconc. Factor (BCF)	6852	pH 4	(1) ACD
Bioconc. Factor (BCF)	14276	pH 7	(1) ACD
Bioconc. Factor (BCF)	14290	pH 8	(1) ACD
Bioconc. Factor (BCF)	14291	pH 10	(1) ACD
Boiling Point (BP)	712.3+-70.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	104.14+-3.0 kJ/mol		(1) ACD
Flash Point (FP)	384.6+-64.2 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	6		(1) ACD
H acceptors (HAC)	8		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	51.6	pH 1	(1) ACD
Koc (KOC)	15714	pH 4	(1) ACD
Koc (KOC)	32737	pH 7	(1) ACD
Koc (KOC)	32769	pH 8	(1) ACD
Koc (KOC)	32772	pH 10	(1) ACD
logD (LOGD)	2.97	pH 1	(1) ACD
logD (LOGD)	5.45	pH 4	(1) ACD
logD (LOGD)	5.77	pH 7	(1) ACD
logD (LOGD)	5.77	pH 8	(1) ACD
logD (LOGD)	5.77	pH 10	(1) ACD
logP (LOGP)	5.771+-0.615		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	584.73		(1) ACD
pKa (PKA)	3.95+-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	3.87E-20 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

=> d 14 1-8 ed

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 03 Feb 2004

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 04 Jan 2002

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 26 Dec 2001

=> d 14 1-8 lc

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CA, CAPLUS, USPATFULL

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED STN Files: CHEMCATS

=> fil hcplus uspatfull chemcats
COST IN U.S. DOLLARS SINCE FILE
TOTAL
SESSION ENTRY
FULL ESTIMATED COST 294.47
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE
TOTAL ENTRY
SESSION CA SUBSCRIBER PRICE -1.32
1.32

FILE 'HCAPLUS' ENTERED AT 10:35:54 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'USPATFULL' ENTERED AT 10:35:54 ON 28 DEC 2004
CA INDEXING COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 American Chemical Society (ACS)

=> d his

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
L1 STRUCTURE UPLOADED
L2 0 S L1 SAM
L3 0 S L1 FAM
L4 8 S L1 FUL
L5 1 S L1 FUL FAM
L6 0 L5 NOT L4

FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004

=> s 14 9 L4

=> d 19 ed

L9 NOT FOUND
The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (>).

=> d 17 1-9 ed

NO VALID FORMATS ENTERED FOR FILE 'USPATFULL'
NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
In a multifile environment, each file must have at least one valid format requested. Refer to file specific help messages or the STNGUIDE file for information on formats available in individual files.

REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

=> d 17 1-9 ibib abs

NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
In a multifile environment, each file must have at least one valid format requested. Refer to file specific help messages or the STNGUIDE file for information on formats available in individual files.

REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):

REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):filedefault

L7 ANSWER 1 OF 9 HCAPLUS COPYRIGHT 2004 ACS on STN

AN 2004:41501 HCPLUS Full-text
DN 140:87744
TI Affinity small molecules for the EPO receptor
IN Olsson, Lennart; Naranda, Tatjana
PA Receptor, Inc., USA
SO PCT Int. Appl., 85 pp.
CODEN: PIXXD2

DT Patent
LA English
FAN.CNT 1
DATE PATENT NO. KIND DATE APPLICATION NO.

PI WO 2004005323 A2 20040115 WO 2003-US21394

20030703 WO 2004005323 A3 20040701

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA,

CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,

LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ,

OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT,

TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW

RW: GH, GM, KE, ES, MW, MZ, SD, SL, TZ, UG, ZM, ZW, AM,

AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,

EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI,

SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN,

TD, TG US 2004171541 A1 20040902 US 2003-613754

20030702 US 2004116346 A1 20040617 US 2003-612885

20030703 PRAI US 2002-393360P P 20020703

US 2002-393361P P 20020703

US 2002-394110P P 20020703

OS MARPAT 140:87744

L7 ANSWER 2 OF 9 USPATFULL on STN
AN 2004:221770 USPATFULL Full-text

TI Affinity small molecules for the EPO receptor
IN Olsson, Lennart, Orinda, CA, UNITED STATES

PI US 2004171541 A1 20040902

AI US 2003-613754 A1 20030702 (10)

PRAI US 2002-393361P 20020703 (60)

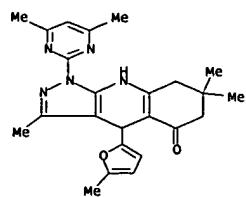
US 2002-393360P 20020703 (60)

US 2002-394110P 20020703 (60)

DT utility

FS APPLICATION

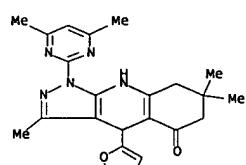
LN.CNT 2046
 INCL INCLM: 514/012.000
 INCLs: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 NCL NCLM: 514/012.000
 NCLs: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 IC [7]
 ICM: A61K038-18
 ICS: A61K031-50; A61K031-519; A61K031-44; A61K031-4745
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.



L7 ANSWER 3 OF 9 USPATFULL on STN
 AN 2004:152124 USPATFULL Full-text
 TI Affinity small molecules for the EPO receptor
 IN Olsson, Lennart, Orinda, CA, UNITED STATES
 Naranda, Tatjana, Mountain View, CA, UNITED STATES
 PI US 2004116346 A1 20040617
 AI US 2003-612885 A1 20030703 (10)
 PRAI US 2002-393361P 20020703 (60)
 US 2002-393360P 20020703 (60)
 US 2002-394110P 20020703 (60)
 DT Utility
 FS APPLICATION
 LN.CNT 2000
 INCL INCLM: 514/012.000
 INCLs: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 NCL NCLM: 514/012.000
 NCLs: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 IC [7]
 ICM: A61K038-17
 ICS: A61K031-519; A61K031-501; A61K031-4745; A61K031-44
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.

L7 ANSWER 4 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:326248 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4266
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-1,4,6,7,8,9-
 hexahydro-
 3,7,7-trimethyl-4-(5-methyl-2-furanyl)-
 CAS Registry No. (RN): 521318-71-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

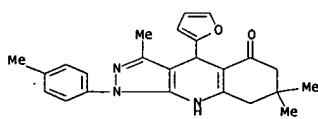
L7 ANSWER 5 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:3262236 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4092
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-
 furanyl)-
 1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-
 1H-pyrazolo[3,4-b]quinolin-5-one



L7 ANSWER 6 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2881171 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-
 CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY

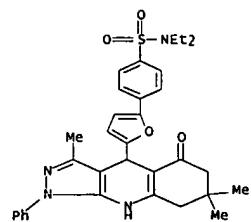
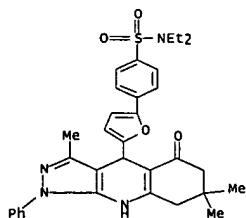
Structure :

CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



L7 ANSWER 7 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2877484 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-
 1H-
 pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

L7 ANSWER 9 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1491648 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-
 1H-
 pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



L7 ANSWER 8 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1507742 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-

>> d 17 1-9 pd
 NO VALID FORMATS ENTERED FOR FILE 'HCAPLUS'
 NO VALID FORMATS ENTERED FOR FILE 'USPATFULL'
 In a multifile environment, each file must have at least one valid
 format requested. Refer to file specific help messages or the
 STNGUIDE file for information on formats available in individual

files.
 REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

>> index biosci
 FILE 'DRUGMON0' ACCESS NOT AUTHORIZED
 COST IN U.S. DOLLARS SINCE FILE
 TOTAL ENTRY
 SESSION 19.21
 FULL ESTIMATED COST
 313.89 SINCE FILE
 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) ENTRY
 TOTAL 0.00 SINCE FILE
 SESSION CA SUBSCRIBER PRICE
 1.32
 1.32
 INDEX 'ADISCT1, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE, AQUALINE, AQUASCI, BIOBUSINESS, BIOCOMMERCE, BIOENG, BIOSIS, BIOTECHABS, BIOTECHDS, BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN, CONFSCI, CROPB, CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON 28 DEC 2004

75 FILES IN THE FILE LIST IN STNINDEX
 Enter SET DETAIL ON to see search term postings or to view search error messages that display as 0* with SET DETAIL OFF.

>> s (epo or (erythropoietin receptor))
 FILE 'ADISCT1'
 68 EPO
 3 EPOS
 71 EPO
 (EPO OR EPOS)
 641 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 642 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 77336 RECEPTOR
 5437 RECEPTORS
 79811 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 3 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'ADISINSIGHT'
 24 EPO
 2 EPOS
 25 EPO
 (EPO OR EPOS)
 57 "ERYTHROPOIETIN"
 15 "ERYTHROPOIETINS"
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'AQUALINE'
 1 EPO
 2 EPOS
 3 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 241 RECEPTOR
 286 RECEPTORS
 395 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'AQUASCI'
 36 EPO
 63 EPOS
 99 EPO
 (EPO OR EPOS)
 10 "ERYTHROPOIETIN"
 5274 "RECEPTOR"
 4864 "RECEPTORS"
 7710 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")
 99 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOBUSINESS'
 387 EPO
 18 EPOS
 404 EPO
 (EPO OR EPOS)
 588 "ERYTHROPOIETIN"
 2 "ERYTHROPOIETINS"
 590 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 6807 "RECEPTOR"
 2441 "RECEPTORS"
 7936 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")
 410 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOCOMMERCE'
 1192 EPO
 3 EPOS
 1197 EPO
 (EPO OR EPOS)
 837 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 837 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 3060 RECEPTOR
 1032 RECEPTORS
 57 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 4128 "RECEPTOR"
 1761 "RECEPTORS"
 4546 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 6 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")
 27 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'ADISNEWS'
 16 EPO
 162 ERYTHROPOIETIN
 5 ERYTHROPOIETINS
 163 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 1962 RECEPTOR
 556 RECEPTORS
 2258 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 3 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 19 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'AGRICOLA'
 35 EPO
 26 EPOS
 60 EPO
 (EPO OR EPOS)
 136 ERYTHROPOIETIN
 8600 RECEPTOR
 9839 RECEPTORS
 13992 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 3 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 63 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'ANABSTR'
 26 EPO
 27 EPOS
 52 EPO
 (EPO OR EPOS)
 72 ERYTHROPOIETIN
 1181 RECEPTOR
 381 RECEPTORS
 1391 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 53 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'ANTE'
 10 EPO
 12 EPOS
 22 EPO
 (EPO OR EPOS)
 7 ERYTHROPOIETIN
 91 RECEPTOR
 73 RECEPTORS
 150 RECEPTOR
 3917 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1197 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOENG'
 180 EPO
 7 EPOS
 183 EPO
 (EPO OR EPOS)
 335 ERYTHROPOIETIN
 9 ERYTHROPOIETINS
 336 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 11378 RECEPTOR
 8432 RECEPTORS
 14157 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 28 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 199 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOSIS'
 5622 EPO
 222 EPOS
 5822 EPO
 (EPO OR EPOS)
 17873 ERYTHROPOIETIN
 64 ERYTHROPOIETINS
 17897 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 651542 RECEPTOR
 323743 RECEPTORS
 780807 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1356 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 6546 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOTECHABS'
 356 EPO
 8 EPOS
 356 EPO
 (EPO OR EPOS)
 981 ERYTHROPOIETIN
 6 ERYTHROPOIETINS
 983 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 16255 RECEPTOR
 4183 RECEPTORS
 17462 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 85 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 407 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOTECHDS'
 356 EPO
 8 EPOS
 356 EPO

(EPO OR EPOS)
 981 ERYTHROPOIETIN
 6 ERYTHROPOETINS
 983 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 16255 RECEPTOR
 4183 RECEPTORS
 17462 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 85 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 407 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'BIOTECHNO'
 2598 EPO
 33 EPOS
 2614 EPO
 (EPO OR EPOS)
 9222 ERYTHROPOIETIN
 29 ERYTHROPOETINS
 9222 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 202158 RECEPTOR
 78475 RECEPTORS
 213334 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 774 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2960 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CABA'
 247 EPO
 47 EPOS
 294 EPO
 (EPO OR EPOS)
 539 ERYTHROPOIETIN
 30692 RECEPTOR
 27815 RECEPTORS
 42614 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 21 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 308 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CANCERLIT'
 3530 EPO
 31 EPOS
 3544 EPO
 (EPO OR EPOS)
 11104 ERYTHROPOIETIN
 30 ERYTHROPOETINS
 11104 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 122617 RECEPTOR
 117416 RECEPTORS
 160576 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 501 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3775 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CONFSCI'
 539 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 87 EPO
 10 EPOS
 97 EPO
 (EPO OR EPOS)
 525 "ERYTHROPOIETIN"
 2 "ERYTHROPOETINS"
 527 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOETINS")
 17313 "RECEPTOR"
 8593 "RECEPTORS"
 25653 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 71 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")
 165 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CROPB'
 0 EPO
 9 EPOS
 9 EPO
 (EPO OR EPOS)
 0 ERYTHROPOIETIN
 160 RECEPTOR
 35 RECEPTORS
 173 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 9 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CROPU'
 55 EPO
 12 EPOS
 67 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 1368 RECEPTOR
 570 RECEPTORS
 1566 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 67 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'DDFB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'DDFU'
 1419 EPO
 3 EPOS
 1421 EPO

FILE 'CAPLUS'
 5521 EPO
 131 EPOS
 5625 EPO
 (EPO OR EPOS)
 11125 ERYTHROPOIETIN
 520 ERYTHROPOETINS
 11155 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 582865 RECEPTOR
 534615 RECEPTORS
 693859 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1237 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 6216 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CEABA-VTB'
 106 EPO
 16 EPOS
 122 EPO
 (EPO OR EPOS)
 191 ERYTHROPOIETIN
 9 ERYTHROPOETINS
 196 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 2128 RECEPTOR
 850 RECEPTORS
 2456 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 4 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 124 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CEN'
 33 EPO
 67 "ERYTHROPOIETIN"
 439 "RECEPTOR"
 381 "RECEPTORS"
 630 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 1 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")
 34 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'CIN'
 533 EPO
 6 EPOS
 536 EPO
 (EPO OR EPOS)
 544 "ERYTHROPOIETIN"
 4 "ERYTHROPOETINS"
 547 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOETINS")
 4037 "RECEPTOR"
 1359 "RECEPTORS"
 4824 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)RECEPTOR")

FILE 'DGENE'
 15416 EPO
 68 EPOS
 15418 EPO
 (EPO OR EPOS)
 18662 ERYTHROPOIETIN
 21 ERYTHROPOETINS
 18680 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOETINS)
 496912 RECEPTOR
 170378 RECEPTORS
 537954 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1551 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 15707 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'DISSABS'
 152 EPO
 51 EPOS
 200 EPO
 (EPO OR EPOS)
 240 ERYTHROPOIETIN
 22949 RECEPTOR
 13830 RECEPTORS
 27515 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 35 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 216 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'DRUGB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))

FILE 'DRUGMONOG2'
 74 EPO
 2 EPOS
 76 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN

0 RECEPTOR
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 76 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'DRUGU'
 1633 EPO
 7 EPOS
 1637 EPO
 (EPO OR EPOS)
 3396 ERYTHROPOETIN
 10 ERYTHROPOETINS
 3396 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 110636 RECEPTOR
 54912 RECEPTORS
 125028 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 64 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 1657 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'EMBAL'
 52 EPO
 2 EPOS
 54 EPO
 (EPO OR EPOS)
 106 ERYTHROPOETIN
 2 ERYTHROPOETINS
 108 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 5008 RECEPTOR
 2397 RECEPTORS
 5899 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 6 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 57 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'EMBASE'
 4482 EPO
 111 EPOS
 4572 EPO
 (EPO OR EPOS)
 18291 "ERYTHROPOETIN"
 51 "ERYTHROPOETINS"
 18292 "ERYTHROPOETIN"
 ("ERYTHROPOETIN" OR "ERYTHROPOETINS")
 714332 "RECEPTOR"
 260226 "RECEPTORS"
 755991 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 1098 ERYTHROPOETIN RECEPTOR
 ("ERYTHROPOETIN(W)RECEPTOR")
 5095 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'ESBIOBASE'
 1925 EPO
 44 EPOS
 1960 EPO
 (EPO OR EPOS)

3664 ERYTHROPOETIN
 11 ERYTHROPOETINS
 3666 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 239241 RECEPTOR
 129772 RECEPTORS
 280062 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 433 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 2150 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'FEDRIP'
 118 EPO
 2 EPOS
 120 EPO
 (EPO OR EPOS)
 243 ERYTHROPOETIN
 16896 RECEPTOR
 9138 RECEPTORS
 19085 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 30 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 139 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'FOMAD'
 1 EPO
 12 EPOS
 13 EPO
 (EPO OR EPOS)
 1 ERYTHROPOETIN
 2 RECEPTOR
 1 RECEPTORS
 2 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 13 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'FOREGE'
 0 EPO
 0 ERYTHROPOETIN
 0 RECEPTOR
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 0 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'FROSTI'
 15 EPO
 15 EPOS
 29 EPO
 (EPO OR EPOS)
 9 ERYTHROPOETIN
 1085 RECEPTOR
 784 RECEPTORS
 1582 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 29 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'FSTA'
 21 EPO
 5 EPOS
 25 EPO
 (EPO OR EPOS)
 1 ERYTHROPOETIN
 1012 RECEPTOR
 434 RECEPTORS
 1297 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 25 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'GENBANK'
 6183 EPO
 6652 "ERYTHROPOETIN"
 201129 "RECEPTOR"
 289 ERYTHROPOETIN RECEPTOR
 ("ERYTHROPOETIN(W)RECEPTOR")
 6457 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'HEALSAFE'
 7 EPO
 19 "ERYTHROPOETIN"
 460 "RECEPTOR"
 311 "RECEPTORS"
 657 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOETIN RECEPTOR
 ("ERYTHROPOETIN(W)RECEPTOR")
 7 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'IFIPAT'
 716 EPO
 11 EPOS
 725 EPO
 (EPO OR EPOS)
 1216 ERYTHROPOETIN
 24 ERYTHROPOETINS
 1235 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 32172 RECEPTOR
 11668 RECEPTORS
 36258 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 68 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 758 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'IMSDRUGNEWS'
 70 EPO
 161 "ERYTHROPOETIN"
 3340 "RECEPTOR"
 836 "RECEPTORS"
 3873 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOETIN RECEPTOR
 ("ERYTHROPOETIN(W)RECEPTOR")
 74 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'KOSMET'
 5 EPO
 1 EPOS
 6 EPO
 (EPO OR EPOS)
 8 ERYTHROPOETIN
 490 RECEPTOR
 555 RECEPTORS
 710 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 800 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'LIFESCI'
 863 EPO
 50 EPOS
 907 EPO
 (EPO OR EPOS)
 1458 "ERYTHROPOETIN"

11 "ERYTHROPOIETINS"
 1458 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 167986 "RECEPTOR"
 153110 "RECEPTORS"
 216519 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 281 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 1047 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'MEDICON'
 4 EPO
 7 EPOS
 11 EPO
 (EPO OR EPOS)
 22 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 23 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 324 RECEPTOR
 571 RECEPTORS
 827 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'MEDLINE'
 4856 EPO
 99 EPOS
 4929 EPO
 (EPO OR EPOS)
 16393 ERYTHROPOIETIN
 53 ERYTHROPOIETINS
 16397 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 479689 RECEPTOR
 514892 RECEPTORS
 673211 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 714 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 5281 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NIOSHTIC'
 15 EPO
 89 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 89 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 1641 RECEPTOR
 1030 RECEPTORS
 2222 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 15 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NTIS'
 46 EPO

64 EPOS
 108 EPO
 (EPO OR EPOS)
 121 ERYTHROPOIETIN
 25 ERYTHROPOIETINS
 122 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 4927 RECEPTOR
 3307 RECEPTORS
 6501 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 108 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NUTRACEUT'
 11 EPO
 0 ERYTHROPOIETIN
 10 RECEPTOR
 13 RECEPTORS
 22 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'OCEAN'
 24 EPO
 51 EPOS
 75 EPO
 (EPO OR EPOS)
 1 "ERYTHROPOIETIN"
 601 "RECEPTOR"
 581 "RECEPTORS"
 950 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 75 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PASCAL'
 2055 EPO
 127 EPOS
 2173 EPO
 (EPO OR EPOS)
 6424 ERYTHROPOIETIN
 22 ERYTHROPOIETINS
 6427 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 304437 RECEPTOR
 102112 RECEPTORS
 327238 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 334 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2342 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PCTGEN'
 0 EPO
 0 ERYTHROPOIETIN
 8549 RECEPTOR

694 RECEPTORS
 9243 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHAR'
 72 EPO
 85 "ERYTHROPOIETIN"
 9931 "RECEPTOR"
 1111 "RECEPTORS"
 10104 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 47 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHARMAL'
 99 EPO
 3 EPOS
 101 EPO
 (EPO OR EPOS)
 202 ERYTHROPOIETIN
 5 ERYTHROPOIETINS
 206 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 1768 RECEPTOR
 465 RECEPTORS
 2031 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 101 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHIC'
 7 EPO
 6 "ERYTHROPOIETIN"
 18 "RECEPTOR"
 5 "RECEPTORS"
 23 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 7 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHIN'
 841 EPO
 22 EPOS
 859 EPO
 (EPO OR EPOS)
 1142 "ERYTHROPOIETIN"
 18 "ERYTHROPOIETINS"
 1153 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 4441 "RECEPTOR"
 2186 "RECEPTORS"
 5782 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 3 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR")

861 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PROMT'
 2713 EPO
 1902 EPOS
 4590 EPO
 (EPO OR EPOS)
 2025 "ERYTHROPOIETIN"
 34 "ERYTHROPOIETINS"
 2050 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 17544 "RECEPTOR"
 9487 "RECEPTORS"
 22972 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 27 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 4612 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PROUDSDR'
 22 EPO
 43 "ERYTHROPOIETIN"
 2 "ERYTHROPOIETINS"
 44 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 36740 "RECEPTOR"
 19875 "RECEPTORS"
 44173 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 4 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)" RECEPTOR")
 22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PS'
 0 EPO
 1 EPOS
 1 EPO
 (EPO OR EPOS)
 0 ERYTHROPOIETIN
 29 RECEPTOR
 1 RECEPTORS
 30 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'RDISCLOSURE'
 21 EPO
 2 EPOS
 23 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 93 RECEPTOR
 20 RECEPTORS
 106 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 23 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'SCISEARCH'

4531 EPO
 222 EPOS
 4734 EPO
 (EPO OR EPOS)
 16279 ERYTHROPOETIN
 46 ERYTHROPOETINS
 16295 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 583687 RECEPTOR
 306298 RECEPTORS
 721251 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1592 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 5851 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'SYNTHLINE'
 1 EPO
 0 "ERYTHROPOETIN"
 798 "RECEPTOR"
 103 "RECEPTORS"
 876 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOETIN RECEPTOR
 ("ERYTHROPOETIN(W)RECEPTOR")
 1 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'TOXCENTER'
 2126 EPO
 34 EPOS
 2145 EPO
 (EPO OR EPOS)
 6597 ERYTHROPOETIN
 72 ERYTHROPOETINS
 6605 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 264727 RECEPTOR
 166955 RECEPTORS
 322116 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 285 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 2274 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'USPATFULL'
 13847 EPO
 184 EPOS
 13982 EPO
 (EPO OR EPOS)
 7989 ERYTHROPOETIN
 1884 ERYTHROPOETINS
 9673 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 104220 RECEPTOR
 75854 RECEPTORS
 118925 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 773 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 14435 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'USPAT2'
 807 EPO
 13 EPOS
 818 EPO
 (EPO OR EPOS)
 477 ERYTHROPOETIN
 30 ERYTHROPOETINS
 494 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 7293 RECEPTOR
 5119 RECEPTORS
 8267 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 32 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 836 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'VETB'
 0 EPO
 1 ERYTHROPOETIN
 77 RECEPTOR
 25 RECEPTORS
 82 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 0 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'VETU'
 23 EPO
 65 ERYTHROPOETIN
 970 RECEPTOR
 641 RECEPTORS
 1302 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 23 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'WATER'
 0 EPO
 2 EPOS
 2 EPO
 (EPO OR EPOS)
 1 ERYTHROPOETIN
 471 RECEPTOR
 265 RECEPTORS
 641 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 2 (EPO OR (ERYTHROPOETIN RECEPTOR))

FILE 'WPIDS'
 580 EPO
 36 EPOS
 612 EPO
 (EPO OR EPOS)
 1446 ERYTHROPOETIN
 16 ERYTHROPOETINS
 1458 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 27 FILE ADISCTI
 27 FILE ADISINSIGHT
 19 FILE ADISNEWS
 63 FILE AGRICOLA
 53 FILE ANABSTR
 22 FILE ANTE
 3 FILE AQUALINE
 99 FILE AQUASCI
 410 FILE BIOPARTNERS
 1197 FILE BIOCOMMERCE
 199 FILE BIOENG
 6546 FILE BIOSIS
 407 FILE BIOTECHABS
 2960 FILE BIOTECHNO
 308 FILE CABA
 3775 FILE CANCERLIT
 6216 FILE CAPLUS
 124 FILE CEABA-VTB
 34 FILE CEN
 539 FILE CIN
 165 FILE CONFSCI
 9 FILE CROPB
 67 FILE CROPU
 3 FILE DDFB
 1441 FILE DDFU
 15707 FILE DGENE
 216 FILE DISSABS
 3 FILE DRUGB
 76 FILE DRUGMONOG2
 1657 FILE DRUGU
 57 FILE EMBAL
 5095 FILE EMBASE
 2150 FILE ESBIOBASE
 139 FILE FEDRIP
 13 FILE FOMAD
 29 FILE FROSTI
 25 FILE FSTA
 6457 FILE GENBANK
 7 FILE HEALSAFE
 758 FILE IFIPAT
 74 FILE IMSDRUGNEWS
 33 FILE IMSPRODUCT
 36 FILE IMSRESEARCH
 800 FILE JICST-EPLUS
 6 FILE KOSMET
 1047 FILE LIFESCI
 11 FILE MEDICONF
 5281 FILE MEDLINE
 15 FILE NIOSHTIC
 108 FILE NTIS

BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN,
 CONFSCI, CROPB, CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON
 28 DEC 2004

SEA (EPO OR (ERYTHROPOETIN RECEPTOR))

73 FILE ADISCTI
 27 FILE ADISINSIGHT
 19 FILE ADISNEWS
 63 FILE AGRICOLA
 53 FILE ANABSTR
 22 FILE ANTE
 3 FILE AQUALINE
 99 FILE AQUASCI
 410 FILE BIOPARTNERS
 1197 FILE BIOCOMMERCE
 199 FILE BIOENG
 6546 FILE BIOSIS
 407 FILE BIOTECHABS
 2960 FILE BIOTECHNO
 308 FILE CABA
 3775 FILE CANCERLIT
 6216 FILE CAPLUS
 124 FILE CEABA-VTB
 34 FILE CEN
 539 FILE CIN
 165 FILE CONFSCI
 9 FILE CROPB
 67 FILE CROPU
 3 FILE DDFB
 1441 FILE DDFU
 15707 FILE DGENE
 216 FILE DISSABS
 3 FILE DRUGB
 76 FILE DRUGMONOG2
 1657 FILE DRUGU
 57 FILE EMBAL
 5095 FILE EMBASE
 2150 FILE ESBIOBASE
 139 FILE FEDRIP
 13 FILE FOMAD
 29 FILE FROSTI
 25 FILE FSTA
 6457 FILE GENBANK
 7 FILE HEALSAFE
 758 FILE IFIPAT
 74 FILE IMSDRUGNEWS
 33 FILE IMSPRODUCT
 36 FILE IMSRESEARCH
 800 FILE JICST-EPLUS
 6 FILE KOSMET
 1047 FILE LIFESCI
 11 FILE MEDICONF
 5281 FILE MEDLINE
 15 FILE NIOSHTIC
 108 FILE NTIS

(ERYTHROPOETIN OR ERYTHROPOETINS)
 41921 RECEPTOR
 13906 RECEPTORS
 46849 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 80 ERYTHROPOETIN
 (ERYTHROPOETIN(W)RECEPTOR)
 659 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'WPIFV'
 10 EPO
 15 ERYTHROPOETIN
 354 RECEPTOR
 89 RECEPTORS
 389 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 2 ERYTHROPOETIN RECEPTOR
 (ERYTHROPOETIN(W)RECEPTOR)
 10 (EPO OR (ERYTHROPOETIN RECEPTOR))
 FILE 'WPINDEX'
 580 EPO
 36 EPOS
 612 EPO
 (EPO OR EPOS)
 1446 ERYTHROPOETIN
 16 ERYTHROPOETINS
 1458 ERYTHROPOETIN
 (ERYTHROPOETIN OR ERYTHROPOETINS)
 41921 RECEPTOR
 13906 RECEPTORS
 46849 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 80 ERYTHROPOETIN
 (ERYTHROPOETIN(W)RECEPTOR)
 659 (EPO OR (ERYTHROPOETIN RECEPTOR))

L8 QUE (EPO OR (ERYTHROPOETIN RECEPTOR))

=> DIS HIST

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004

L1 STRUCTURE UPLOADED
 L2 0 S L1 SAM
 L3 0 S L1 FAM
 L4 8 S L1 FUL
 L5 1 S L1 FUL FAM
 L6 0 LS NOT 4

FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28
 DEC 2004

L7 9 S L4

INDEX 'ADISCTI, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE,
 AQUALINE,
 AQUASCI, BIOPARTNERS, BIOCOMMERCE, BIOENG, BIOSIS, BIOTECHABS,
 BIOTECHDS,

11 FILE NUTRACEUT
75 FILE OCEAN
2342 FILE PASCAL
73 FILE PHAR
101 FILE PHARMAML
7 FILE PHIC
861 FILE PHIN
4612 FILE PROMT
22 FILE PROUSDDR
1 FILE PS
23 FILE RDISCLOSURE
5851 FILE SCISEARCH
1 FILE SYNTHLINE
2274 FILE TOXCENTER
14435 FILE USPATFULL
836 FILE USPAT2
23 FILE VETU
2 FILE WATER
659 FILE WPIDS
10 FILE WPIFV
659 FILE WPINDEX
L8 QUE (EPO OR (ERYTHROPOIETIN RECEPTOR))

=>

---Logging off of STN---

=>
Executing the logoff script...

=> LOG Y
COST IN U.S. DOLLARS SINCE FILE
TOTAL ENTRY

SESSION
FULL ESTIMATED COST 11.97
325.86

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE
TOTAL ENTRY

SESSION
CA SUBSCRIBER PRICE 0.00
1.32

STN INTERNATIONAL LOGOFF AT 10:50:37 ON 28 DEC 2004